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INFORMATION DISCLOSURE CITATION (Use several sheets if necessary)	ATTY. DOCKET NO. PC10697A	SERIAL NO. 09/804,176
	APPLICANT S. Ekins	
	FILING DATE March 12, 2001	GROUP 1646 /631

U.S. PATENT DOCUMENTS

EXAMINER INITIAL	DOCUMENT NUMBER	DATE	NAME	CLASS	SUBCLASS	FILING DATE IF APPROPRIATE

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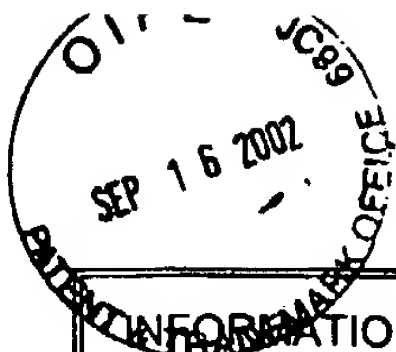
FOREIGN PATENT DOCUMENTS

DOCUMENT NUMBER	DATE	COUNTRY	CLASS	SUBCLASS	TRANSLATION
					YES NO

OTHER DOCUMENTS (Including Author, Title, Date, Pertinent Pages, Etc.)

JB	Gunter R. Strobl et al, "Development of a Pharmacophore for Inhibition of Human Liver Cytochrome P-450 2D6: Molecular Modeling and Inhibition Studies", J. Med. Chem. 1993, 96, 1136-1145.
JB	Dafang Wu et al, "Interactions of Amphetamine Analogs with Human Liver CYP2D6", Elsevier Science Inc., Biochemical Pharmacology, Vol. 53, pp. 1605-1612, 1997.
JB	Yoshihiko Nishibata et al, "Automatic Creation of Drug Candidate Structures Based on Receptor Structure. Starting Point for Artificial Lead Generation", Pergamon Press plc, Tetrahedron Vol. 47, No. 43, pp. 8985-8990, 1991.
JB	Sean Ekins et al, "Three and Four Dimensional-Quantitative Structure Activity Relationship (3D/4D-QSAR) Analyses of CYP2D6 Inhibitors", Pharmacogenetics 1999, 9:477-489.
JB	Jonathan Greene et al, "Chemical Function Queries for 3D Database Search", J. Chem. Inf. Comput. Sci. 1994, 34, 1297-1308.
JB	Scott J. Weiner et al, "A New Force Field for Molecular Mechanical Simulation of Nucleic Acids and Proteins", J. Am. Chem. Soc. 1984, 106, 765-784.

JB. Bruce 29 November 2001



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-	JB		P. J. Goodford et al, "A Computational Procedure for Determining Energetically Favorable Binding Sites on Biologically Important Macromolecules", J. Med. Chem. 1985, 28, 849-857.			
-	JB		Andrew Miranker et al, "Functionality Maps of Binding Sites: A Multiple Copy Simultaneous Search Method", Wiley-Liss, Inc., PROTEINS: Structure, Function and Genetics 11:29-34 (1991).			
-	JB		Irwin D. Kuntz et al, "A Geometric Approach to Macromolecule-Ligand Interactions", Academic Press Inc., J. Mol. Biol. (1982) 161, 269-288.			
-	JB		Sean Ekins et al, "Three-Dimensional Quantitative Structure Activity Relationship Analyses of Substrates for CYP2B6", The Journal of Pharmacology and Experimental Therapeutics, Vol. 288, No. 1, pp. 21-29. (1999)			
-	JB		Erik Skjelbo et al, "Inhibitors of Imipramine Metabolism by Human Liver Microsomes", Br. J. clin. Pharmac. (1992), 34, 256-261.			
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